Trying 3106016892...Open

Welcome to STN International! Enter x:x
LOGINID:ssspta1611bxv
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

Welcome to STN International Web Page URLs for STN Seminar Schedule - N. America NEWS The CA Lexicon available in the CAPLUS and CA files NEWS Dec 17 NEWS Feb 06 Engineering Information Encompass files have new names TOXLINE no longer being updated NEWS Feb 16 Apr 23 Search Derwent WPINDEX by chemical structure NEWS Apr 23 PRE-1967 REFERENCES NOW SEARCHABLE IN CAPLUS AND CA NEWS NEWS May 07 DGENE Reload NEWS Jun 20 Published patent applications (A1) are now in USPATFULL 8 NEWS 9 JUL 13 New SDI alert frequency now available in Derwent's DWPI and DPCI In-process records and more frequent updates now in NEWS 10 Aug 23 MEDLINE PAGE IMAGES FOR 1947-1966 RECORDS IN CAPLUS AND CA NEWS 11 Aug 23 Adis Newsletters (ADISNEWS) now available on STN NEWS 12 Aug 23 NEWS 13 Sep 17 IMSworld Pharmaceutical Company Directory name change to PHARMASEARCH Korean abstracts now included in Derwent World Patents NEWS 14 Oct 09 Index NEWS 15 Oct 09 Number of Derwent World Patents Index updates increased NEWS 16 Oct 15 Calculated properties now in the REGISTRY/ZREGISTRY File NEWS 17 Oct 22 Over 1 million reactions added to CASREACT NEWS 18 Oct 22 DGENE GETSIM has been improved NEWS 19 Oct 29 AAASD no longer available NEWS 20 Nov 19 New Search Capabilities USPATFULL and USPAT2 NEWS 21 Nov 19 TOXCENTER(SM) - new toxicology file now available on STN NEWS 22 Nov 29 COPPERLIT now available on STN NEWS 23 Nov 29 DWPI revisions to NTIS and US Provisional Numbers NEWS 24 Nov 30 Files VETU and VETB to have open access NEWS 25 Dec 10 WPINDEX/WPIDS/WPIX New and Revised Manual Codes for 2002 NEWS 26 Dec 10 DGENE BLAST Homology Search NEWS EXPRESS August 15 CURRENT WINDOWS VERSION IS V6.0c, CURRENT MACINTOSH VERSION IS V6.0 (ENG) AND V6.0J (JP), AND CURRENT DISCOVER FILE IS DATED 07 AUGUST 2001 NEWS HOURS STN Operating Hours Plus Help Desk Availability NEWS INTER General Internet Information NEWS LOGIN Welcome Banner and News Items Direct Dial and Telecommunication Network Access to STN NEWS PHONE CAS World Wide Web Site (general information) NEWS WWW

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FILE 'HOME' ENTERED AT 16:59:42 ON 13 DEC 2001

=> file reg COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.15 0.15

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 16:59:58 ON 13 DEC 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2001 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 11 DEC 2001 HIGHEST RN 374745-93-4 DICTIONARY FILE UPDATES: 11 DEC 2001 HIGHEST RN 374745-93-4

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=>
Uploading C:\STNEXP4\QUERIES\09685053.str

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> d l1

L1 HAS NO ANSWERS

L1

STR

G1 O,S,N,Cy,Ak,[@1]

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam SAMPLE SEARCH INITIATED 17:00:34 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 10351 TO ITERATE

9.7% PROCESSED 1000 ITERATIONS INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED) SEARCH TIME: 00.00.01 16 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**

BATCH **COMPLETE**

PROJECTED ITERATIONS: 200934 TO 213106 PROJECTED ANSWERS: 2540 TO 4084

L3 16 SEA SSS SAM L1

=> d scan

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN Acetamide, N-[4-[(1-methylethyl)amino]-1,3,5-triazin-2-yl]- (9CI)

MF C8 H13 N5 O

09/679,150

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):15

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN Benzeneacetamide,
3-[[4-[(5-phenyl-1H-pyrazol-3-yl)amino]-1,3,5-triazin-2yl]amino]- (9CI)
MF C20 H18 N8 O

*** FRAGMENT DIAGRAM IS INCOMPLETE ***

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN Benzenesulfonamide,
4-methoxy-N-[4-[(1-phenylethyl)amino]-1,3,5-triazin-2yl]- (9CI)
MF C18 H19 N5 O3 S

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN 1,3,5-Triazine-2,4-diamine, N-heptyl-N'-hexadecyl- (9CI)
MF C26 H51 N5

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1H-Benzimidazol-2-amine, N-(3-pyridinylmethyl)-1-[4-[(3,4,5-trimethoxyphenyl)amino]-1,3,5-triazin-2-yl]- (9CI)

MF C25 H24 N8 O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN 1,3,5-Triazine-2,4-diamine,
N-methyl-N-[2-[4-(1-methylethoxy)phenyl]ethyl](9CI)

MF C15 H21 N5 O

CI COM

$$\begin{array}{c|c} & \text{Me} \\ & \text{N} & \text{N} \\ & \text{N} & \text{CH}_2 - \text{CH}_2 \\ & \text{OPr-i} \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN Benzeneacetic acid, 2-[[4-(2-ethoxyphenoxy)-1,3,5-triazin-2-yl]oxy]-

.alpha.-(methoxymethylene)-, methyl ester, (E)- (9CI) C22 H21 N3 O6 MF

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

REGISTRY COPYRIGHT 2001 ACS L3 16 ANSWERS

1,3,5-Triazine-2,4-diamine, N-[1,1'-biphenyl]-4-yl-N'-(3,4,5-IN trimethoxyphenyl) - (9CI)

C24 H23 N5 O3 MF

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3

16 ANSWERS REGISTRY COPYRIGHT 2001 ACS
1,3,5-Triazin-2-amine, N-[4-(4-morpholinyl)phenyl]-4-[2-(phenylamino)-3-IN pyridinyl] - (9CI)

C24 H23 N7 O MF

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN Ethenamine, N, N-dimethyl-2-(1,3,5-triazin-2-yl)- (9CI)

MF C7 H10 N4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1H-Isoindole-1,3(2H)-dione, 2-(4-amino-1,3,5-triazin-2-yl)- (9CI)

MF C11 H7 N5 O2

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N-(3,4-dimethoxyphenyl)-N'-(4-phenoxyphenyl)(9CI)

MF C23 H21 N5 O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS IN 1,3,5-Triazin-2-amine, N,N-diphenyl- (9CI)

MF C15 H12 N4

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2-acetonitrile, 4-methyl-.alpha.-[2-[4-[(5-methyl-2-furanyl)oxy]phenoxy]-1-oxopropyl]- (9CI)

MF C20 H18 N4 O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N,N'-bis(3-chlorophenyl)-N,N'-dimethyl- (9CI)

MF C17 H15 Cl2 N5

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 16 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N'-ethyl-N,N-di-2-propenyl- (9CI)

MF C11 H17 N5

$$\begin{array}{c|c} \operatorname{CH}_2-\operatorname{CH} = \operatorname{CH}_2 \\ | & \\ \operatorname{EtNH} & \operatorname{N} - \operatorname{CH}_2 - \operatorname{CH} = \operatorname{CH}_2 \\ | & \\ \operatorname{N} & \operatorname{N} \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

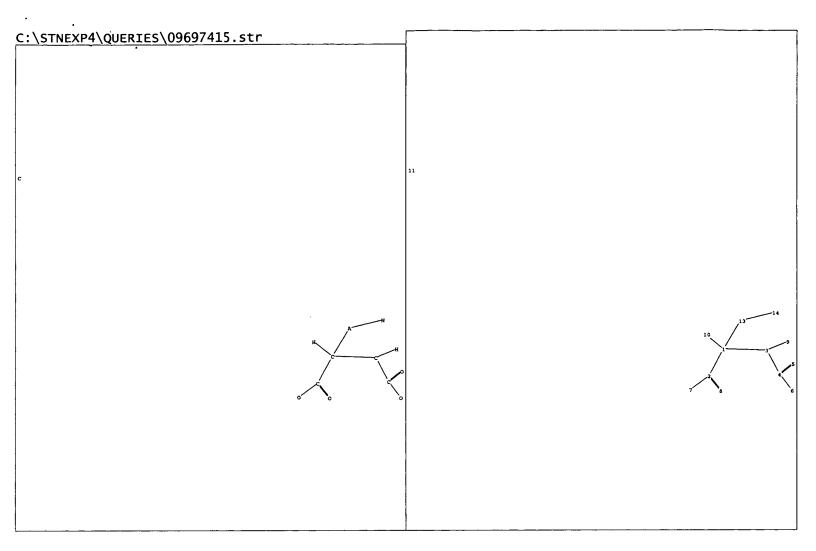
=> log y

COST IN U.S. DOLLARS SINCE FILE TOTAL

FULL ESTIMATED COST ENTRY SESSION 0.62 0.77

STN INTERNATIONAL LOGOFF AT 17:01:14 ON 13 DEC 2001

Connection closed by remote host



chain nodes :
 1 2 3 4 5 6 7 8 9 10 11 13 14
chain bonds :
 1-2 1-3 1-10 1-13 2-7 2-8 3-4 3-9 4-5 4-6 13-14
exact/norm bonds :
 1-13 2-7 2-8 4-5 4-6 13-14
exact bonds :
 1-2 1-3 1-10 3-4 3-9

Match level:
1:CLASS 2:CLASS 3:CLASS 4:CLASS 5:CLASS 6:CLASS 7:CLASS 8:CLASS 9:CLASS 10:CLASS 11;CLASS 13:CLASS 14:CLASS

6016892...Open

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LOGINID:ssspta1611bxv
PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2

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FILE 'HOME' ENTERED AT 16:13:22 ON 10 DEC 2001

=> file reg

COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.15 0.15

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 16:13:31 ON 10 DEC 2001 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2001 American Chemical Society (ACS)

STRUCTURE FILE UPDATES: 9 DEC 2001 HIGHEST RN 374591-02-3 DICTIONARY FILE UPDATES: 9 DEC 2001 HIGHEST RN 374591-02-3

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=>Testing the current file.... screen

ENTER SCREEN EXPRESSION OR (END):end

=> Uploading C:\STNEXP4\QUERIES\09685053.str

L1 STRUCTURE UPLOADED

=> que L1

L2 QUE L1

=> d l1

L1 HAS NO ANSWERS

L1 STR



G1 O, S, N, Cy, Ak, [@1]

Structure attributes must be viewed using STN Express query preparation.

=> s l1 sss sam SAMPLE SEARCH INITIATED 16:14:09 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 10338 TO ITERATE

1000 ITERATIONS 9.7% PROCESSED INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED) SEARCH TIME: 00.00.02

14 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**

PROJECTED ITERATIONS: 200678 TO 212842

2173 TO PROJECTED ANSWERS: 3615

14 SEA SSS SAM L1

=> d scan

14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

Benzeneacetamide,

3-[[4-[(5-phenyl-1H-pyrazol-3-yl)amino]-1,3,5-triazin-2-

yl]amino]- (9CI)

C20 H18 N8 O

*** FRAGMENT DIAGRAM IS INCOMPLETE ***

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):13

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN Acetamide, N-[4-[(1-methylethyl)amino]-1,3,5-triazin-2-yl]- (9CI)

MF C8 H13 N5 O

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS IN 1,3,5-Triazine-2,4-diamine, N-methyl-N-[2-[4-(1-methylethoxy)phenyl]ethyl]-(9CI)

MF C15 H21 N5 O

CI COM

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1H-Benzimidazol-2-amine, N-(3-pyridinylmethyl)-1-[4-[(3,4,5-

trimethoxyphenyl)amino]-1,3,5-triazin-2-yl]- (9CI)

MF C25 H24 N8 O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N-heptyl-N'-hexadecyl- (9CI)

MF C26 H51 N5

Me⁻ (CH₂)₆-NH
$$\stackrel{\circ}{}$$
N NH- (CH₂)₁₅-Me

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazin-2-amine, N-[4-(4-morpholinyl)phenyl]-4-[2-(phenylamino)-3pyridinyl]- (9CI)

MF C24 H23 N7 O

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N-[1,1'-biphenyl]-4-yl-N'-(3,4,5-

trimethoxyphenyl) - (9CI)

MF C24 H23 N5 O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN Benzeneacetic acid, 2-[[4-(2-ethoxyphenoxy)-1,3,5-triazin-2-yl]oxy].alpha.-(methoxymethylene)-, methyl ester, (E)- (9CI)

MF C22 H21 N3 O6

Double bond geometry as shown.

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N-(3,4-dimethoxyphenyl)-N'-(4-phenoxyphenyl)(9CI)

MF C23 H21 N5 O3

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1H-Isoindole-1,3(2H)-dione, 2-(4-amino-1,3,5-triazin-2-yl)- (9CI)

MF C11 H7 N5 O2

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N'-ethyl-N,N-di-2-propenyl- (9CI)

MF C11 H17 N5

$$\begin{array}{c|c} \text{CH}_2-\text{CH} & \text{CH}_2\\ \\ \vdash & \text{N-CH}_2-\text{CH} & \text{CH}_2\\ \\ N & N & \end{array}$$

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2,4-diamine, N,N'-bis(3-chlorophenyl)-N,N'-dimethyl- (9CI)

MF C17 H15 Cl2 N5

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS

IN 1,3,5-Triazine-2-acetonitrile, 4-methyl-.alpha.-[2-[4-[(5-methyl-2-furanyl)oxy]phenoxy]-1-oxopropyl]- (9CI)

MF C20 H18 N4 O4

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L3 14 ANSWERS REGISTRY COPYRIGHT 2001 ACS
IN Benzenesulfonamide,
4-methoxy-N-[4-[(1-phenylethyl)amino]-1,3,5-triazin-2yl]- (9CI)
MF C18 H19 N5 O3 S

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> log y COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION 0.62 0.77

FULL ESTIMATED COST

::::j{:::